

# ***Interactive comment on “Application of random forest regression to the calculation of gas-phase chemistry within the GEOS-Chem chemistry model v10” by Christoph A. Keller and Mat J. Evans***

**Christoph A. Keller and Mat J. Evans**

christoph.a.keller@nasa.gov

Received and published: 16 January 2019

The comment was uploaded in the form of a supplement:  
<https://www.geosci-model-dev-discuss.net/gmd-2018-229/gmd-2018-229-AC2-supplement.pdf>

Interactive comment on Geosci. Model Dev. Discuss., <https://doi.org/10.5194/gmd-2018-229>, 2018.

Printer-friendly version

Discussion paper

